



# Full wwPDB X-ray Structure Validation Report ⓘ

Nov 9, 2020 – 10:48 AM EST

PDB ID : 6V8G  
Title : GltPh mutant - Y204L A345V V366A  
Authors : Boudker, O.; Huysmans, G.H.M.  
Deposited on : 2019-12-11  
Resolution : 3.38 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at [validation@mail.wwpdb.org](mailto:validation@mail.wwpdb.org)

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

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The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.8.5 (274361), CSD as541be (2020)  
Xtriage (Phenix) : 1.13  
EDS : 2.14.6  
buster-report : 1.1.7 (2018)  
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)  
Refmac : 5.8.0158  
CCP4 : 7.0.044 (Gargrove)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.14.6

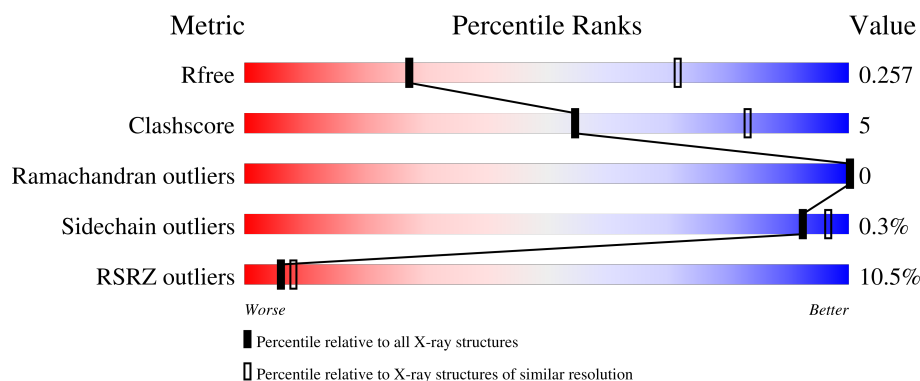
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 3.38 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
$R_{free}$	130704	1691 (3.46-3.30)
Clashscore	141614	1762 (3.46-3.30)
Ramachandran outliers	138981	1732 (3.46-3.30)
Sidechain outliers	138945	1731 (3.46-3.30)
RSRZ outliers	127900	1635 (3.46-3.30)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions  $\leq 5\%$ . The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	422	
1	B	422	
1	C	422	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	ASP	A	503	-	-	X	-
3	ASP	B	503	-	-	X	-
3	ASP	C	503	-	-	X	-

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 8679 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called Glutamate transporter homolog.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	393	Total	C	N	O	S	0	0	0
			2881	1895	463	506	17			
1	B	393	Total	C	N	O	S	0	0	0
			2881	1895	463	506	17			
1	C	393	Total	C	N	O	S	0	0	0
			2881	1895	463	506	17			

There are 42 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	37	HIS	ASP	conflict	UNP O59010
A	40	HIS	LYS	conflict	UNP O59010
A	125	HIS	LYS	conflict	UNP O59010
A	132	HIS	LYS	conflict	UNP O59010
A	204	LEU	TYR	engineered mutation	UNP O59010
A	223	HIS	LYS	conflict	UNP O59010
A	264	HIS	LYS	conflict	UNP O59010
A	345	VAL	ALA	engineered mutation	UNP O59010
A	366	ALA	VAL	engineered mutation	UNP O59010
A	368	HIS	GLU	conflict	UNP O59010
A	418	THR	GLU	conflict	UNP O59010
A	420	VAL	-	expression tag	UNP O59010
A	421	PRO	-	expression tag	UNP O59010
A	422	ARG	-	expression tag	UNP O59010
B	37	HIS	ASP	conflict	UNP O59010
B	40	HIS	LYS	conflict	UNP O59010
B	125	HIS	LYS	conflict	UNP O59010
B	132	HIS	LYS	conflict	UNP O59010
B	204	LEU	TYR	engineered mutation	UNP O59010
B	223	HIS	LYS	conflict	UNP O59010
B	264	HIS	LYS	conflict	UNP O59010
B	345	VAL	ALA	engineered mutation	UNP O59010
B	366	ALA	VAL	engineered mutation	UNP O59010

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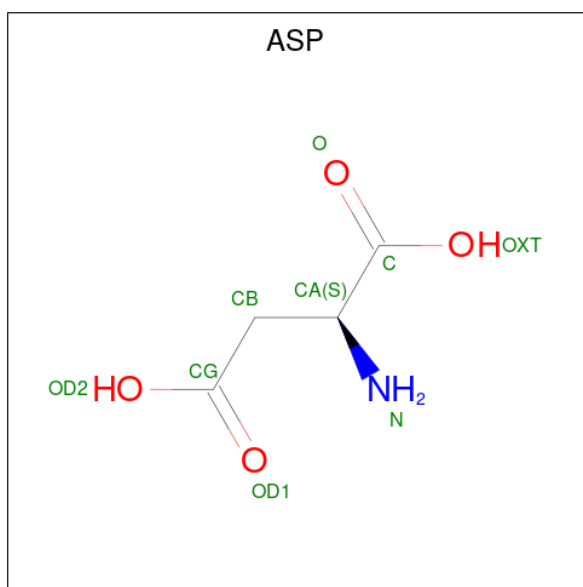
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Chain	Residue	Modelled	Actual	Comment	Reference
B	368	HIS	GLU	conflict	UNP O59010
B	418	THR	GLU	conflict	UNP O59010
B	420	VAL	-	expression tag	UNP O59010
B	421	PRO	-	expression tag	UNP O59010
B	422	ARG	-	expression tag	UNP O59010
C	37	HIS	ASP	conflict	UNP O59010
C	40	HIS	LYS	conflict	UNP O59010
C	125	HIS	LYS	conflict	UNP O59010
C	132	HIS	LYS	conflict	UNP O59010
C	204	LEU	TYR	engineered mutation	UNP O59010
C	223	HIS	LYS	conflict	UNP O59010
C	264	HIS	LYS	conflict	UNP O59010
C	345	VAL	ALA	engineered mutation	UNP O59010
C	366	ALA	VAL	engineered mutation	UNP O59010
C	368	HIS	GLU	conflict	UNP O59010
C	418	THR	GLU	conflict	UNP O59010
C	420	VAL	-	expression tag	UNP O59010
C	421	PRO	-	expression tag	UNP O59010
C	422	ARG	-	expression tag	UNP O59010

- Molecule 2 is SODIUM ION (three-letter code: NA) (formula: Na) (labeled as "Ligand of Interest" by author).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	B	3	Total Na 3 3	0	0
2	A	3	Total Na 3 3	0	0
2	C	3	Total Na 3 3	0	0

- Molecule 3 is ASPARTIC ACID (three-letter code: ASP) (formula: C<sub>4</sub>H<sub>7</sub>NO<sub>4</sub>) (labeled as "Ligand of Interest" by author).

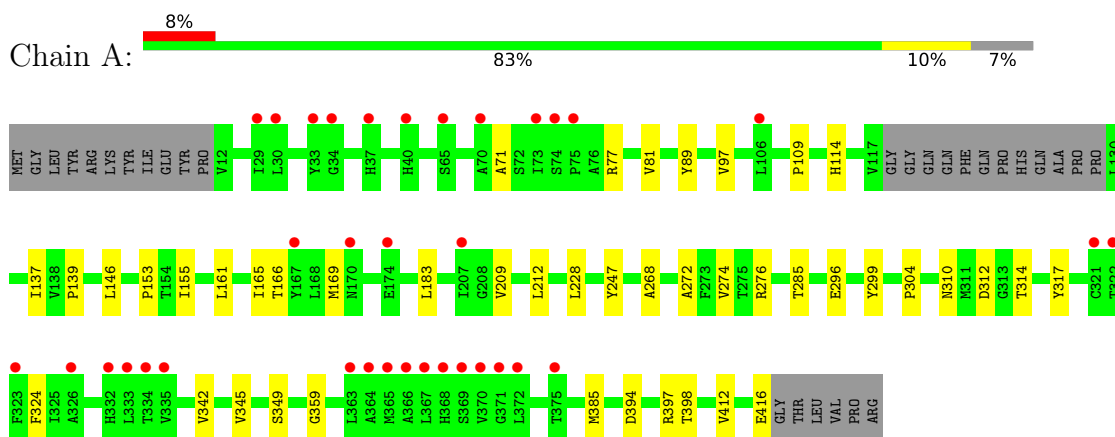


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			9	4	1	4		
3	B	1	Total	C	N	O	0	0
			9	4	1	4		
3	C	1	Total	C	N	O	0	0
			9	4	1	4		

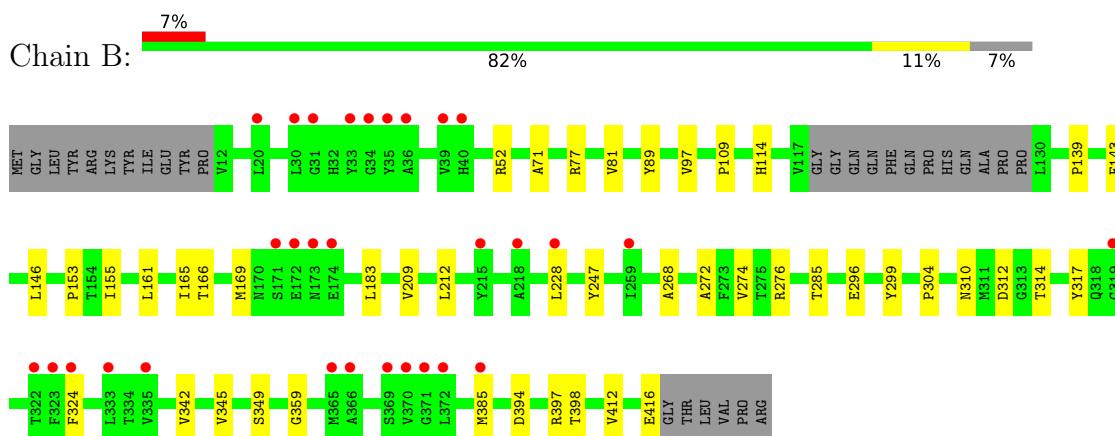
### 3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ( $RSRZ > 2$ ). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

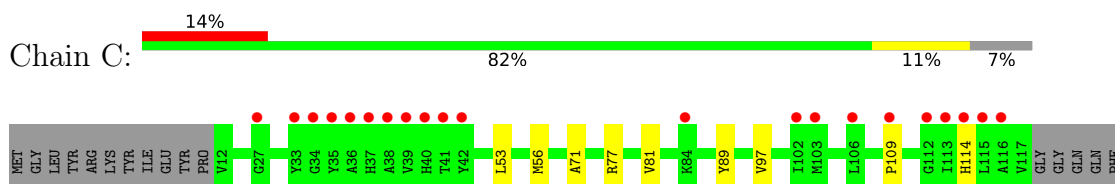
- Molecule 1: Glutamate transporter homolog

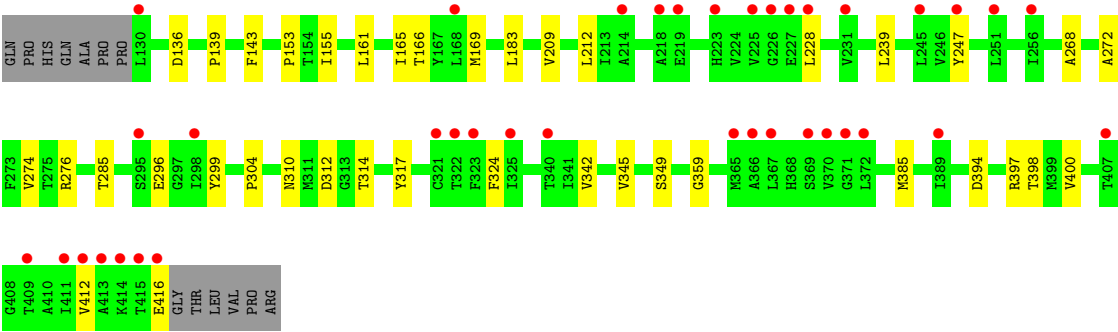


- Molecule 1: Glutamate transporter homolog



- Molecule 1: Glutamate transporter homolog







## 4 Data and refinement statistics

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	109.42Å 109.42Å 305.23Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	47.38 – 3.38 48.18 – 3.38	Depositor EDS
% Data completeness (in resolution range)	99.6 (47.38-3.38) 92.4 (48.18-3.38)	Depositor EDS
$R_{merge}$	0.12	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.62 (at 3.40Å)	Xtriage
Refinement program	PHENIX 1.16_3549	Depositor
R, $R_{free}$	0.232 , 0.257 0.232 , 0.257	Depositor DCC
$R_{free}$ test set	1400 reflections (4.57%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	105.8	Xtriage
Anisotropy	0.504	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.30 , 82.0	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.028 for -h,-k,l	Xtriage
$F_o, F_c$ correlation	0.89	EDS
Total number of atoms	8679	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	132.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.46% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical values of  $\langle |L| \rangle$ ,  $\langle L^2 \rangle$  for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

## 5 Model quality

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section:  
NA

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with  $|Z| > 5$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z  >5	RMSZ	# Z  >5
1	A	0.25	0/2930	0.39	0/3996
1	B	0.25	0/2930	0.39	0/3996
1	C	0.26	0/2930	0.39	0/3996
All	All	0.25	0/8790	0.39	0/11988

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

### 5.2 Too-close contacts

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2881	0	3049	30	0
1	B	2881	0	3049	32	0
1	C	2881	0	3049	34	0
2	A	3	0	0	0	0
2	B	3	0	0	0	0
2	C	3	0	0	0	0
3	A	9	0	3	4	0
3	B	9	0	3	5	0
3	C	9	0	3	5	0
All	All	8679	0	9156	93	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 5.

All (93) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:155:ILE:HD11	1:B:304:PRO:HB2	1.70	0.72
1:A:155:ILE:HD11	1:A:304:PRO:HB2	1.71	0.72
1:C:155:ILE:HD11	1:C:304:PRO:HB2	1.72	0.72
1:C:268:ALA:HB1	1:C:285:THR:HG22	1.79	0.64
1:B:268:ALA:HB1	1:B:285:THR:HG22	1.80	0.63
1:A:268:ALA:HB1	1:A:285:THR:HG22	1.79	0.63
1:C:209:VAL:HA	1:C:274:VAL:HG11	1.83	0.60
1:B:209:VAL:HA	1:B:274:VAL:HG11	1.84	0.59
1:C:398:THR:HG1	3:C:503:ASP:N	2.00	0.59
1:B:398:THR:HG1	3:B:503:ASP:N	2.01	0.59
1:A:209:VAL:HA	1:A:274:VAL:HG11	1.83	0.59
1:C:81:VAL:HG13	1:C:412:VAL:HG11	1.86	0.58
1:A:398:THR:HG1	3:A:503:ASP:N	2.01	0.58
1:B:81:VAL:HG13	1:B:412:VAL:HG11	1.84	0.58
1:A:209:VAL:HG13	1:A:274:VAL:HG21	1.85	0.58
1:A:81:VAL:HG13	1:A:412:VAL:HG11	1.85	0.57
1:C:209:VAL:HG13	1:C:274:VAL:HG21	1.85	0.57
1:B:209:VAL:HG13	1:B:274:VAL:HG21	1.85	0.57
1:C:394:ASP:OD1	3:C:503:ASP:N	2.39	0.56
1:A:394:ASP:OD1	3:A:503:ASP:N	2.39	0.56
1:C:97:VAL:HG11	1:C:342:VAL:HA	1.88	0.55
1:B:97:VAL:HG11	1:B:342:VAL:HA	1.88	0.55
1:A:228:LEU:HD21	1:A:385:MET:HB3	1.89	0.55
1:C:359:GLY:N	3:C:503:ASP:OD2	2.37	0.55
1:A:97:VAL:HG11	1:A:342:VAL:HA	1.88	0.55
1:C:314:THR:HA	1:C:397:ARG:HD2	1.89	0.54
1:A:314:THR:HA	1:A:397:ARG:HD2	1.89	0.54
1:B:314:THR:HA	1:B:397:ARG:HD2	1.90	0.54
1:B:394:ASP:OD1	3:B:503:ASP:N	2.40	0.53
1:C:228:LEU:HD21	1:C:385:MET:HB3	1.90	0.53
1:B:276:ARG:HD2	1:B:394:ASP:HB3	1.91	0.52
1:A:359:GLY:N	3:A:503:ASP:OD2	2.37	0.52
1:B:228:LEU:HD21	1:B:385:MET:HB3	1.90	0.52
1:C:166:THR:HA	1:C:169:MET:HG2	1.91	0.52
1:A:276:ARG:HD2	1:A:394:ASP:HB3	1.91	0.51
1:B:345:VAL:O	1:B:349:SER:N	2.41	0.51
1:C:345:VAL:O	1:C:349:SER:N	2.41	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:166:THR:HA	1:B:169:MET:HG2	1.93	0.51
1:B:359:GLY:N	3:B:503:ASP:OD2	2.37	0.51
1:A:166:THR:HA	1:A:169:MET:HG2	1.92	0.51
1:C:139:PRO:HB3	1:C:153:PRO:HB3	1.93	0.51
1:C:276:ARG:HD2	1:C:394:ASP:HB3	1.92	0.51
1:A:139:PRO:HB3	1:A:153:PRO:HB3	1.93	0.50
1:A:272:ALA:HB1	1:A:398:THR:HG22	1.94	0.50
1:C:247:TYR:OH	1:C:312:ASP:OD2	2.25	0.50
1:B:247:TYR:OH	1:B:312:ASP:OD2	2.25	0.49
1:A:345:VAL:O	1:A:349:SER:N	2.41	0.49
1:B:139:PRO:HB3	1:B:153:PRO:HB3	1.94	0.49
1:C:272:ALA:HB1	1:C:398:THR:HG22	1.95	0.49
1:A:71:ALA:HB1	1:A:166:THR:HG23	1.95	0.48
1:A:146:LEU:HB3	1:B:143:PHE:CD2	2.48	0.48
1:B:272:ALA:HB1	1:B:398:THR:HG22	1.95	0.48
1:C:109:PRO:HB2	1:C:324:PHE:HD1	1.79	0.47
1:C:71:ALA:HB1	1:C:166:THR:HG23	1.97	0.47
1:A:247:TYR:OH	1:A:312:ASP:OD2	2.24	0.47
1:A:109:PRO:HB2	1:A:324:PHE:HD1	1.80	0.45
1:B:146:LEU:HB3	1:C:143:PHE:CD2	2.50	0.45
1:C:161:LEU:O	1:C:165:ILE:HG12	2.16	0.45
1:B:109:PRO:HB2	1:B:324:PHE:HD1	1.80	0.45
1:B:71:ALA:HB1	1:B:166:THR:HG23	1.97	0.45
1:B:212:LEU:HB2	1:B:274:VAL:HG13	1.98	0.45
1:C:212:LEU:HB2	1:C:274:VAL:HG13	1.97	0.45
1:A:161:LEU:O	1:A:165:ILE:HG12	2.17	0.44
1:B:161:LEU:O	1:B:165:ILE:HG12	2.16	0.44
1:A:212:LEU:HB2	1:A:274:VAL:HG13	1.98	0.44
1:B:77:ARG:HH11	1:B:416:GLU:HB3	1.82	0.44
1:C:239:LEU:HD22	1:C:400:VAL:HG21	2.00	0.44
1:A:77:ARG:HH11	1:A:416:GLU:HB3	1.82	0.44
1:C:77:ARG:HH11	1:C:416:GLU:HB3	1.83	0.44
1:C:71:ALA:O	1:C:166:THR:HG21	2.18	0.44
1:B:397:ARG:NE	3:B:503:ASP:OD1	2.50	0.43
1:B:52:ARG:NH2	1:C:136:ASP:HA	2.34	0.43
1:C:397:ARG:NE	3:C:503:ASP:OD1	2.49	0.43
1:A:71:ALA:O	1:A:166:THR:HG21	2.19	0.43
1:A:296:GLU:HA	1:A:299:TYR:CE2	2.54	0.42
1:A:165:ILE:HD11	1:A:183:LEU:HD23	2.01	0.42
1:B:71:ALA:O	1:B:166:THR:HG21	2.19	0.42
1:B:165:ILE:HD11	1:B:183:LEU:HD23	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:296:GLU:HA	1:B:299:TYR:CE2	2.55	0.42
1:C:165:ILE:HD11	1:C:183:LEU:HD23	2.01	0.42
1:C:296:GLU:HA	1:C:299:TYR:CE2	2.55	0.42
1:C:317:TYR:CG	1:C:397:ARG:HD3	2.55	0.42
1:B:89:TYR:CD1	1:B:310:ASN:HB2	2.55	0.41
1:B:398:THR:OG1	3:B:503:ASP:N	2.53	0.41
1:B:317:TYR:CG	1:B:397:ARG:HD3	2.56	0.41
1:A:397:ARG:NE	3:A:503:ASP:OD1	2.50	0.41
1:C:89:TYR:CD1	1:C:310:ASN:HB2	2.55	0.41
1:C:398:THR:OG1	3:C:503:ASP:N	2.53	0.41
1:A:89:TYR:CD1	1:A:310:ASN:HB2	2.56	0.41
1:A:317:TYR:CG	1:A:397:ARG:HD3	2.56	0.41
1:C:53:LEU:HA	1:C:56:MET:HE2	2.03	0.41
1:C:212:LEU:HD12	1:C:274:VAL:HG12	2.04	0.40
1:A:137:ILE:O	1:A:153:PRO:HA	2.22	0.40

There are no symmetry-related clashes.

## 5.3 Torsion angles ⓘ

### 5.3.1 Protein backbone ⓘ

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	389/422 (92%)	383 (98%)	6 (2%)	0	100	100
1	B	389/422 (92%)	382 (98%)	7 (2%)	0	100	100
1	C	389/422 (92%)	382 (98%)	7 (2%)	0	100	100
All	All	1167/1266 (92%)	1147 (98%)	20 (2%)	0	100	100

There are no Ramachandran outliers to report.

### 5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	301/330 (91%)	300 (100%)	1 (0%)	92	96
1	B	301/330 (91%)	300 (100%)	1 (0%)	92	96
1	C	301/330 (91%)	300 (100%)	1 (0%)	92	96
All	All	903/990 (91%)	900 (100%)	3 (0%)	92	96

All (3) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	114	HIS
1	B	114	HIS
1	C	114	HIS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

## 5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

## 5.5 Carbohydrates ⓘ

There are no monosaccharides in this entry.

## 5.6 Ligand geometry ⓘ

Of 12 ligands modelled in this entry, 9 are monoatomic - leaving 3 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with  $|Z| > 2$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z  > 2$	Counts	RMSZ	$\# Z  > 2$
3	ASP	A	503	-	2,8,8	0.31	0	1,10,10	0.32	0
3	ASP	C	503	-	2,8,8	0.30	0	1,10,10	0.36	0
3	ASP	B	503	-	2,8,8	0.31	0	1,10,10	0.36	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	ASP	A	503	-	-	0/2/8/8	-
3	ASP	C	503	-	-	0/2/8/8	-
3	ASP	B	503	-	-	0/2/8/8	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

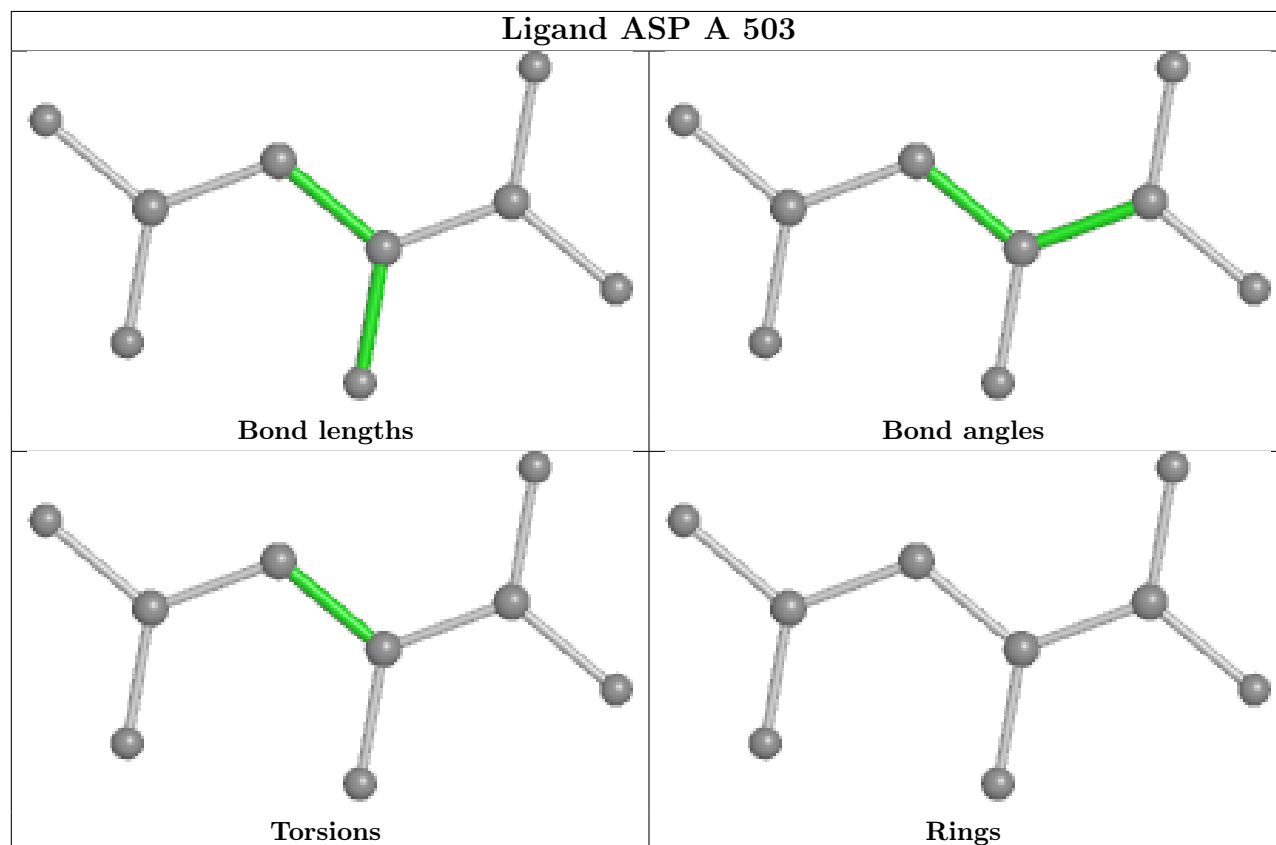
There are no ring outliers.

3 monomers are involved in 14 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	503	ASP	4	0
3	C	503	ASP	5	0
3	B	503	ASP	5	0

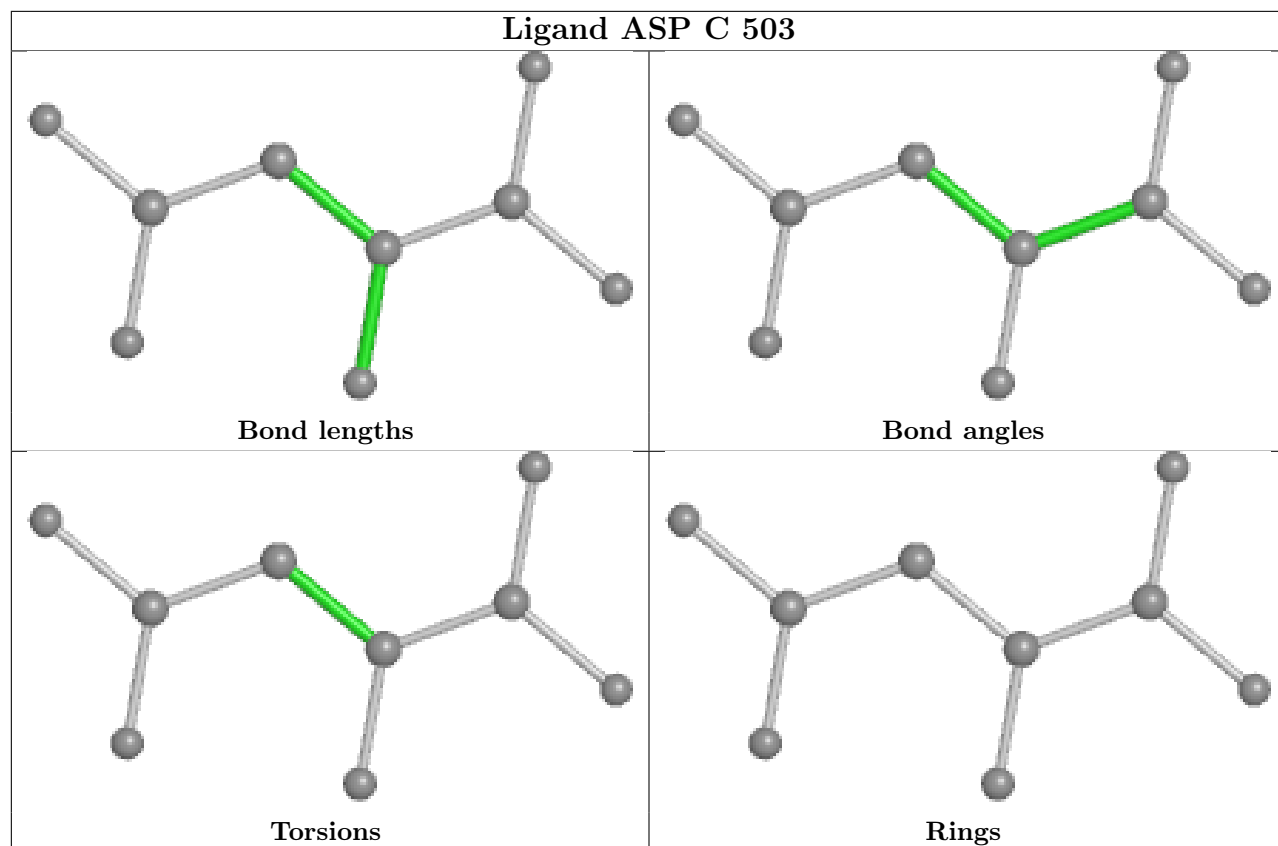
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight  $> 250$  and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring

in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.

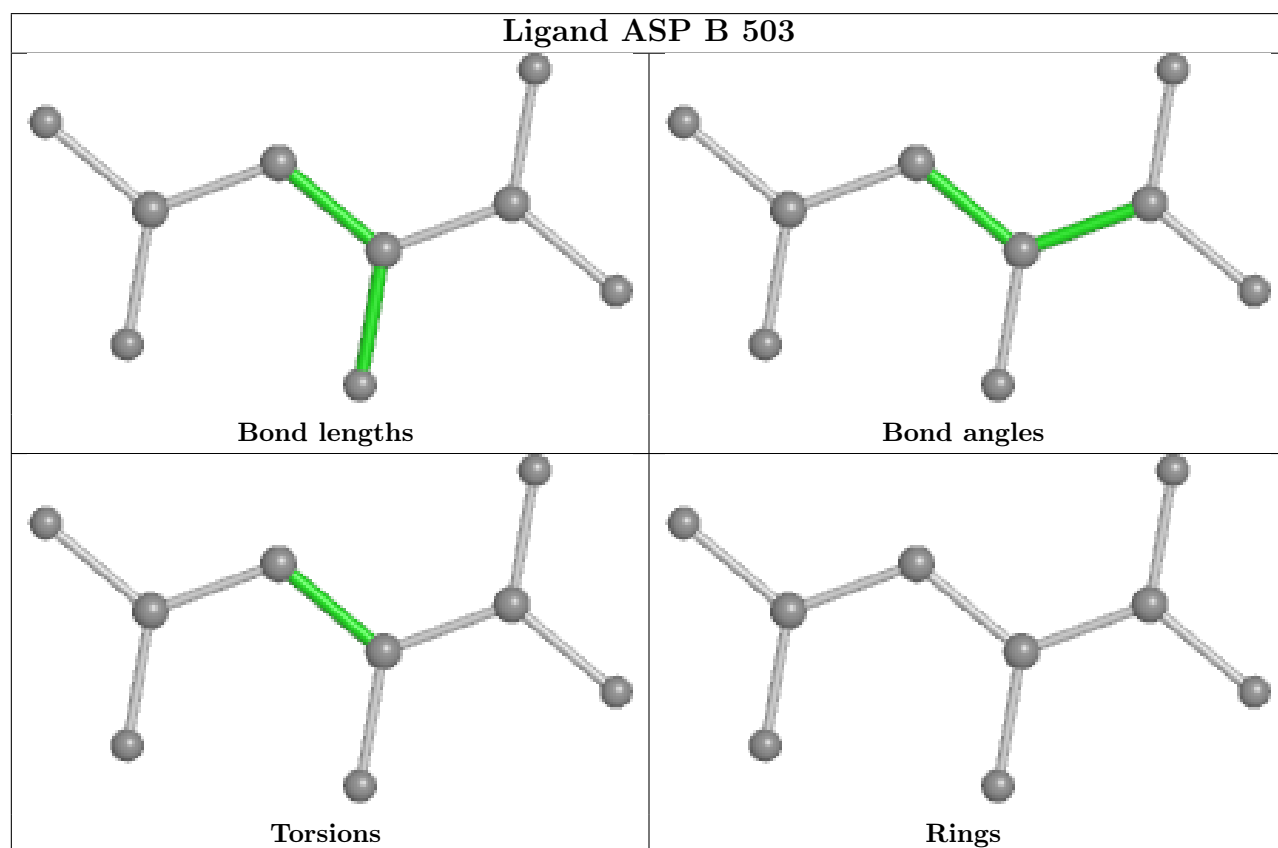




## Ligand ASP C 503



## Ligand ASP B 503



## 5.7 Other polymers [i](#)

There are no such residues in this entry.

## 5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

## 6 Fit of model and data

### 6.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95<sup>th</sup> percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å <sup>2</sup> )	Q<0.9
1	A	393/422 (93%)	0.45	35 (8%) <b>9</b> <b>12</b>	77, 127, 196, 269	0
1	B	393/422 (93%)	0.40	30 (7%) <b>13</b> <b>17</b>	74, 127, 195, 269	0
1	C	393/422 (93%)	0.61	59 (15%) <b>2</b> <b>3</b>	76, 134, 197, 275	0
All	All	1179/1266 (93%)	0.49	124 (10%) <b>6</b> <b>8</b>	74, 129, 197, 275	0

All (124) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	370	VAL	9.8
1	C	369	SER	9.6
1	C	371	GLY	7.9
1	A	367	LEU	6.9
1	C	114	HIS	6.7
1	C	35	TYR	6.3
1	C	372	LEU	6.3
1	B	172	GLU	6.0
1	C	112	GLY	5.8
1	C	36	ALA	5.6
1	C	367	LEU	5.5
1	B	35	TYR	5.4
1	A	333	LEU	4.9
1	A	366	ALA	4.8
1	A	174	GLU	4.5
1	A	372	LEU	4.5
1	C	228	LEU	4.5
1	B	30	LEU	4.4
1	C	412	VAL	4.2
1	A	73	ILE	4.2
1	B	36	ALA	4.1
1	A	371	GLY	4.0
1	B	33	TYR	3.9

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Mol	Chain	Res	Type	RSRZ
1	C	39	VAL	3.8
1	C	366	ALA	3.8
1	B	333	LEU	3.7
1	C	34	GLY	3.7
1	C	219	GLU	3.6
1	A	369	SER	3.6
1	C	37	HIS	3.5
1	C	340	THR	3.5
1	C	322	THR	3.5
1	B	371	GLY	3.5
1	B	369	SER	3.5
1	B	215	TYR	3.4
1	A	323	PHE	3.3
1	B	171	SER	3.3
1	C	298	ILE	3.3
1	C	414	LYS	3.2
1	A	364	ALA	3.2
1	A	106	LEU	3.2
1	C	413	ALA	3.1
1	A	370	VAL	3.1
1	C	218	ALA	3.1
1	A	322	THR	3.1
1	C	416	GLU	3.1
1	C	295	SER	3.0
1	B	31	GLY	3.0
1	A	74	SER	3.0
1	C	227	GLU	3.0
1	B	173	ASN	2.9
1	C	42	TYR	2.9
1	C	106	LEU	2.9
1	A	335	VAL	2.9
1	C	223	HIS	2.9
1	A	70	ALA	2.9
1	A	37	HIS	2.8
1	C	84	LYS	2.8
1	B	366	ALA	2.8
1	C	41	THR	2.8
1	A	33	TYR	2.7
1	A	207	ILE	2.7
1	B	34	GLY	2.7
1	C	116	ALA	2.7
1	C	115	LEU	2.7

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Mol	Chain	Res	Type	RSRZ
1	A	29	ILE	2.7
1	B	322	THR	2.7
1	C	102	ILE	2.6
1	C	38	ALA	2.6
1	C	323	PHE	2.6
1	B	385	MET	2.6
1	C	109	PRO	2.6
1	C	415	THR	2.6
1	B	259	ILE	2.6
1	C	256	ILE	2.6
1	B	335	VAL	2.6
1	C	247	TYR	2.6
1	B	174	GLU	2.6
1	C	113	ILE	2.5
1	A	34	GLY	2.5
1	B	319	GLY	2.5
1	A	326	ALA	2.5
1	C	251	LEU	2.5
1	A	365	MET	2.5
1	B	218	ALA	2.5
1	C	27	GLY	2.5
1	B	323	PHE	2.4
1	A	363	LEU	2.4
1	C	389	ILE	2.4
1	B	228	LEU	2.4
1	B	370	VAL	2.4
1	A	170	ASN	2.4
1	A	332	HIS	2.3
1	C	40	HIS	2.3
1	C	409	THR	2.3
1	A	75	PRO	2.3
1	C	225	VAL	2.3
1	B	20	LEU	2.3
1	B	39	VAL	2.3
1	A	167	TYR	2.3
1	A	321	CYS	2.3
1	C	411	ILE	2.3
1	B	40	HIS	2.2
1	B	372	LEU	2.2
1	A	368	HIS	2.2
1	A	334	THR	2.2
1	C	130	LEU	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	40	HIS	2.2
1	C	103	MET	2.2
1	C	321	CYS	2.2
1	A	375	THR	2.1
1	A	65	SER	2.1
1	B	324	PHE	2.1
1	C	33	TYR	2.1
1	C	214	ALA	2.1
1	C	325	ILE	2.1
1	C	168	LEU	2.1
1	C	407	THR	2.0
1	C	245	LEU	2.0
1	C	365	MET	2.0
1	C	231	VAL	2.0
1	A	30	LEU	2.0
1	C	226	GLY	2.0
1	B	365	MET	2.0

## 6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

## 6.3 Carbohydrates [i](#)

There are no monosaccharides in this entry.

## 6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å <sup>2</sup> )	Q<0.9
3	ASP	B	503	9/9	0.85	0.26	166,171,179,198	0
3	ASP	C	503	9/9	0.86	0.32	180,182,185,202	0
3	ASP	A	503	9/9	0.90	0.28	167,176,182,202	0
2	NA	C	501	1/1	0.92	0.51	92,92,92,92	0
2	NA	C	502	1/1	0.93	0.18	86,86,86,86	0
2	NA	B	504	1/1	0.96	0.10	118,118,118,118	0

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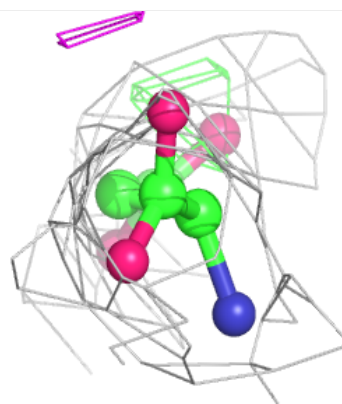
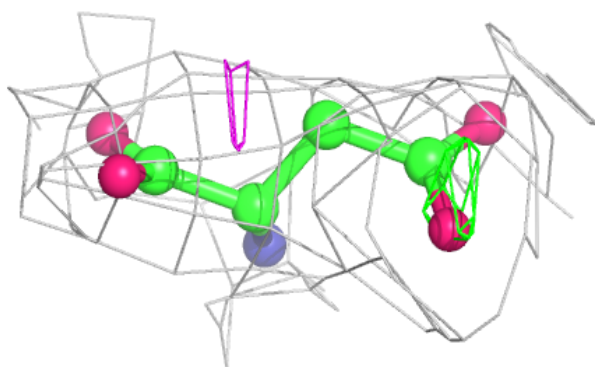
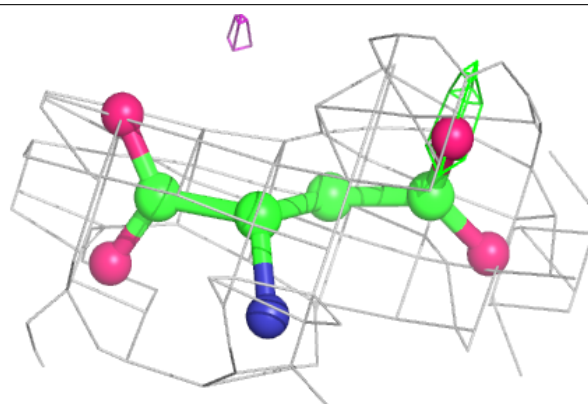
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
2	NA	A	504	1/1	0.97	0.19	99,99,99,99	0
2	NA	A	501	1/1	0.97	0.23	89,89,89,89	0
2	NA	B	501	1/1	0.97	0.41	102,102,102,102	0
2	NA	A	502	1/1	0.97	0.15	82,82,82,82	0
2	NA	B	502	1/1	0.97	0.11	75,75,75,75	0
2	NA	C	504	1/1	0.98	0.10	117,117,117,117	0

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

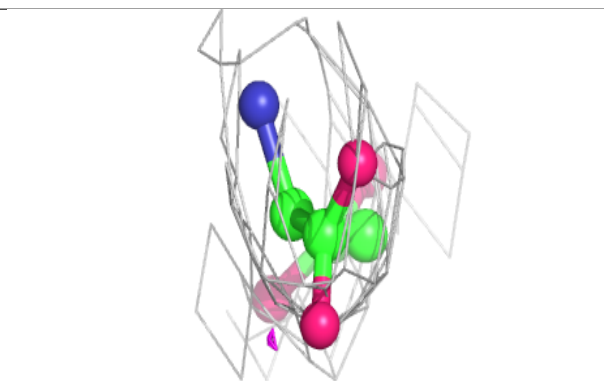
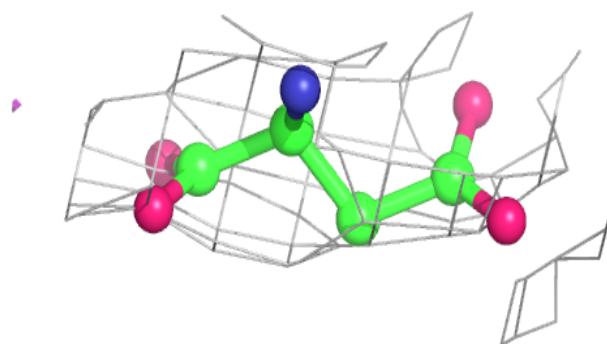
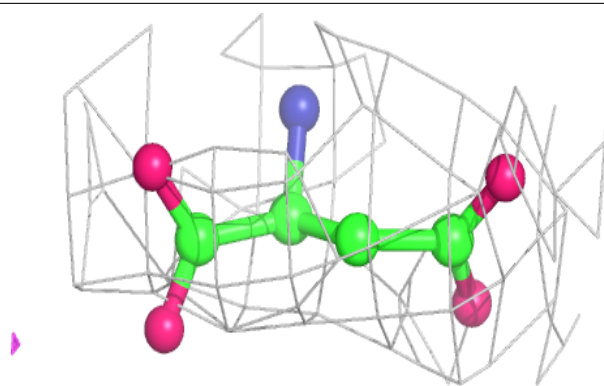
**Electron density around ASP B 503:**

2mF<sub>o</sub>-DF<sub>c</sub> (at 0.7 rmsd) in gray  
mF<sub>o</sub>-DF<sub>c</sub> (at 3 rmsd) in purple (negative)  
and green (positive)

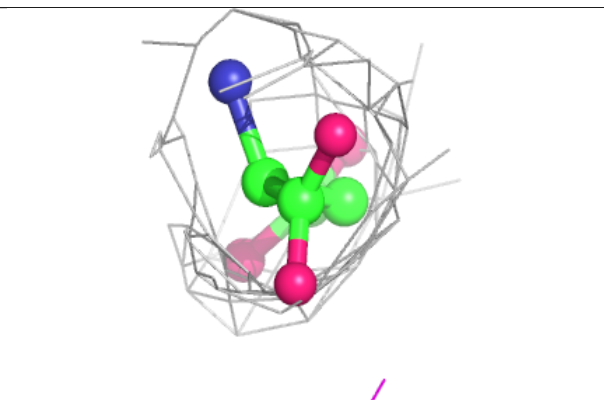
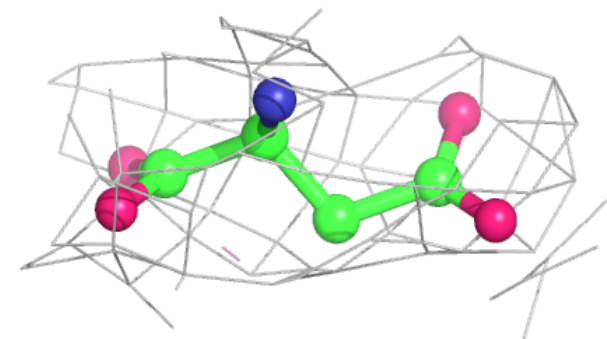
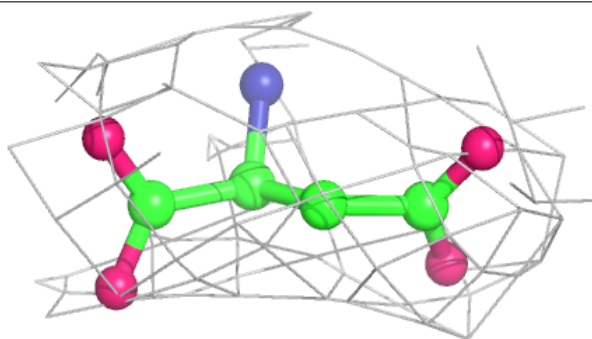


**Electron density around ASP C 503:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

**Electron density around ASP A 503:**

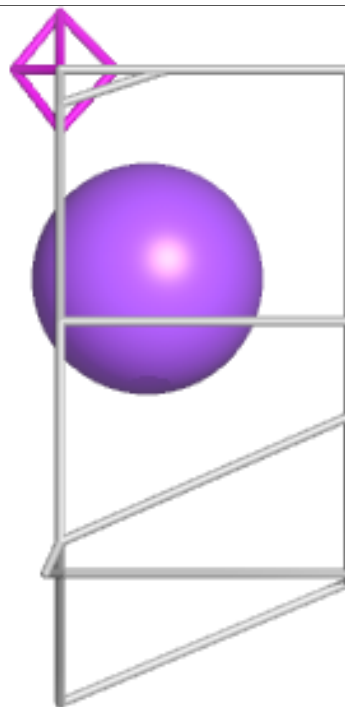
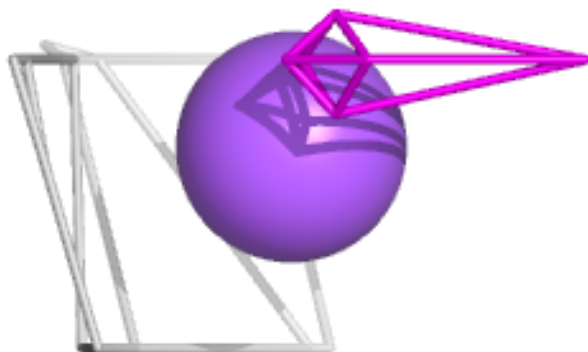
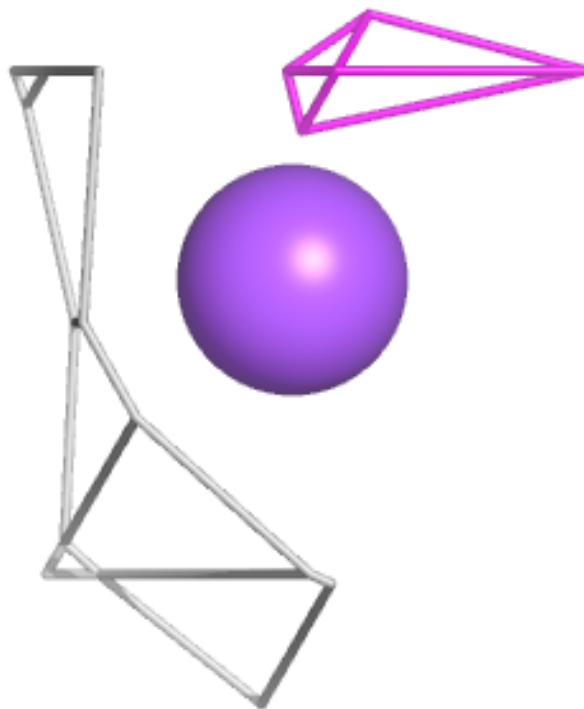
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)





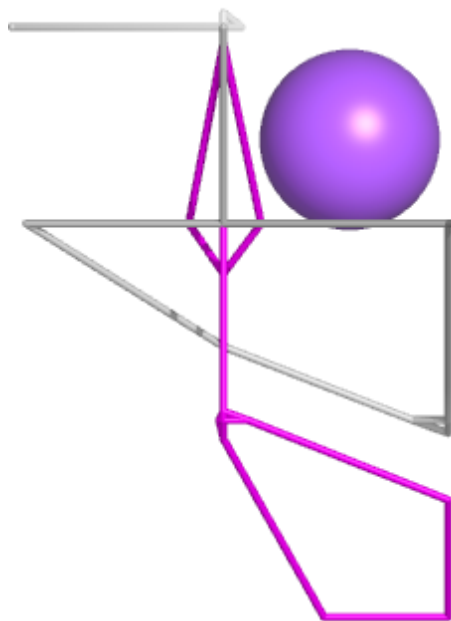
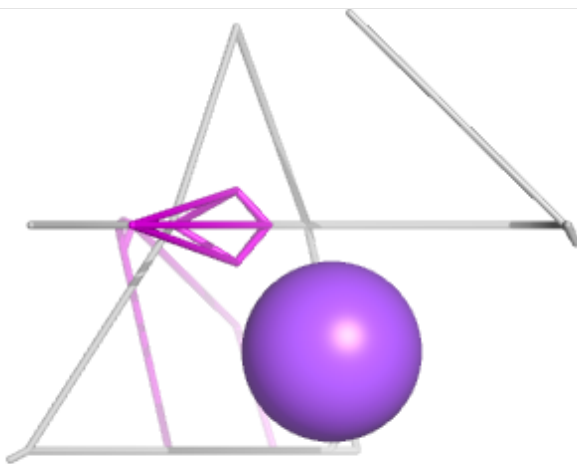
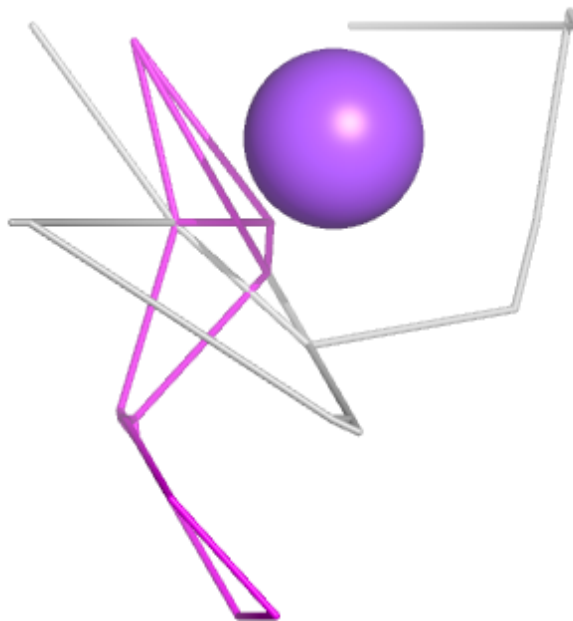
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 $mF_o - DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)



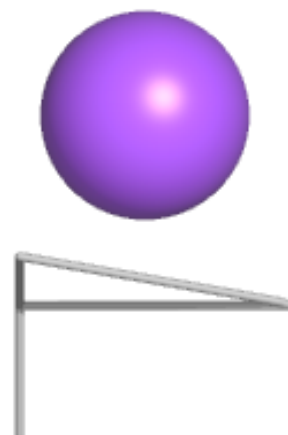
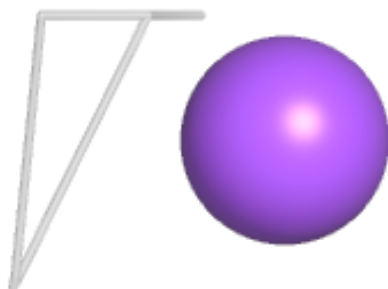
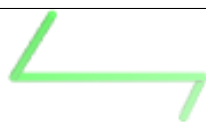
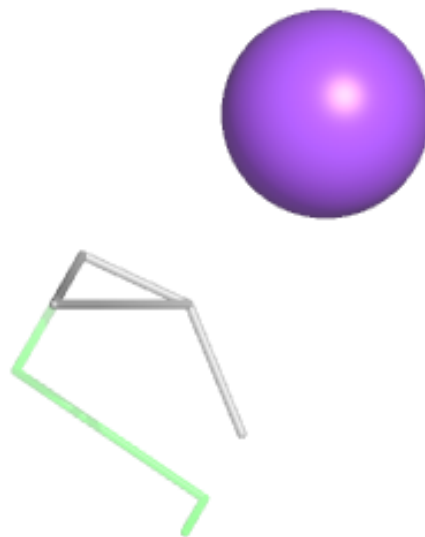
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and green (positive)



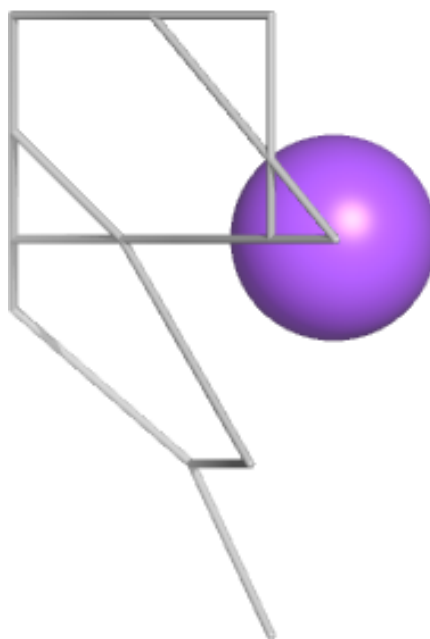
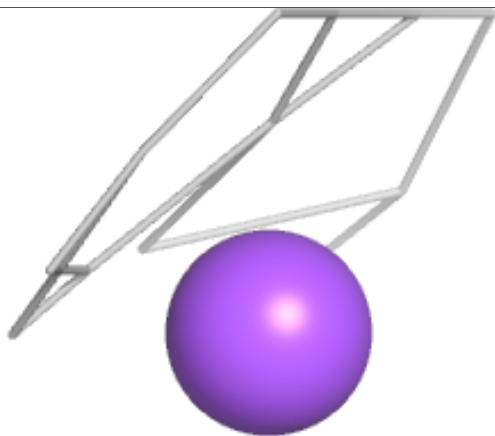
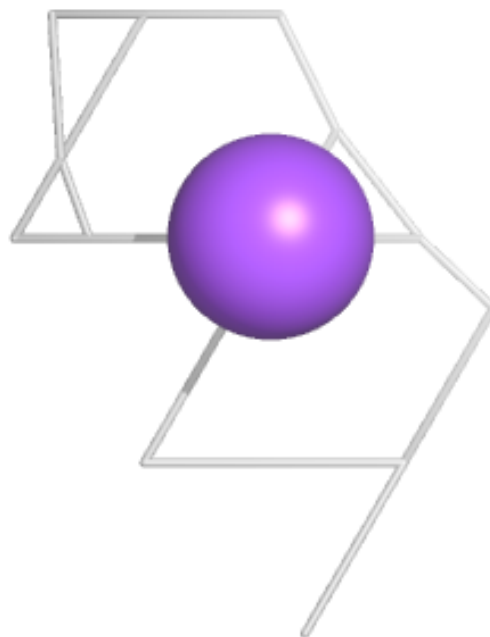
**Electron density around NA B 504:**

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and green (positive)



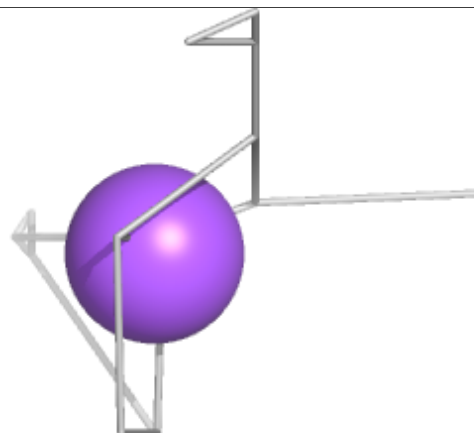
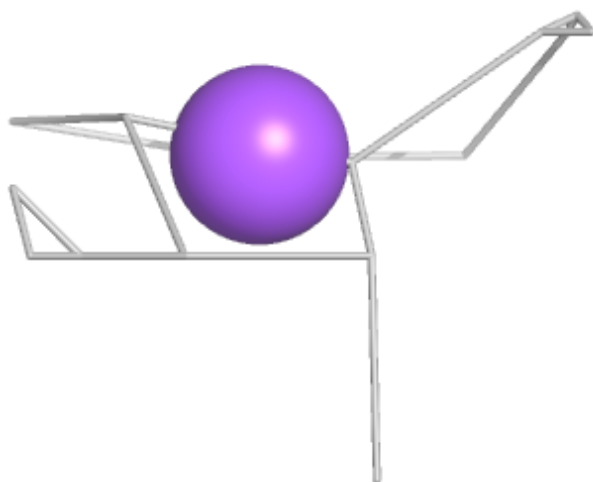
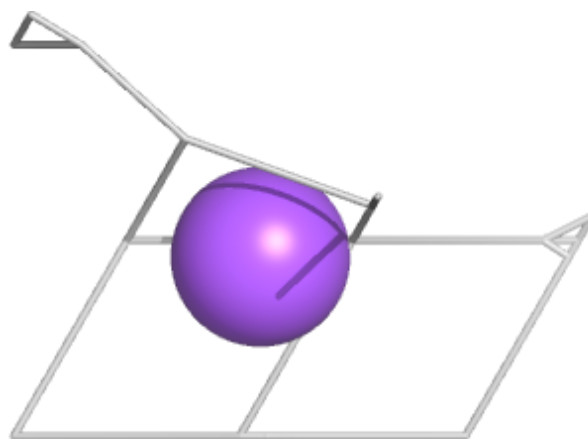
**Electron density around NA A 504:**

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and green (positive)



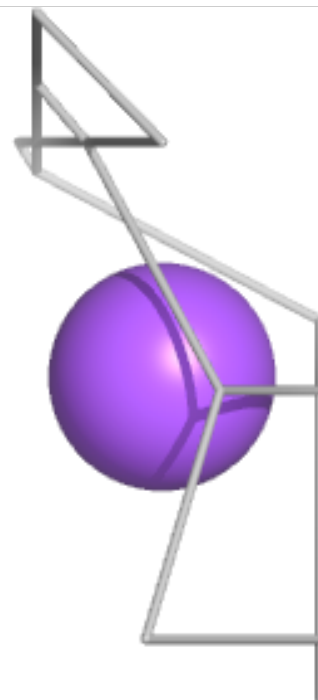
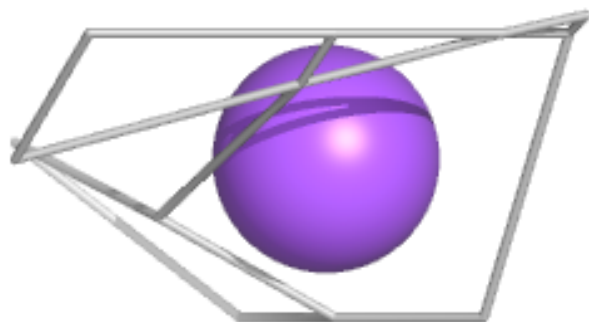
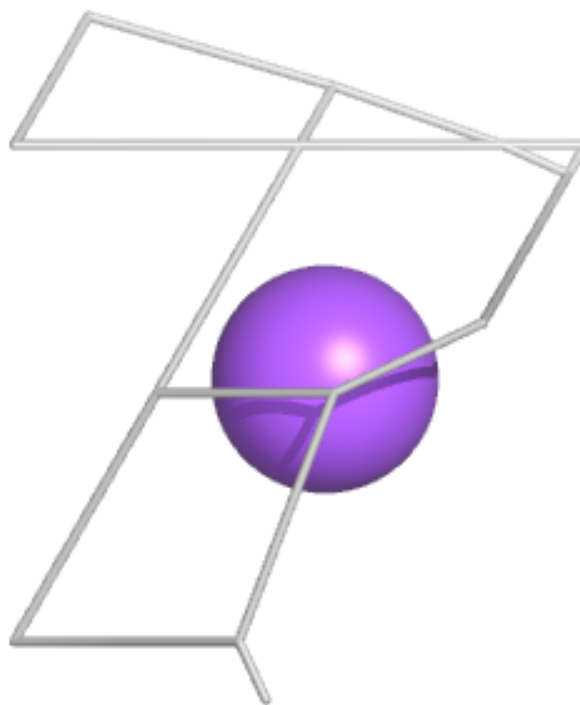
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and green (positive)



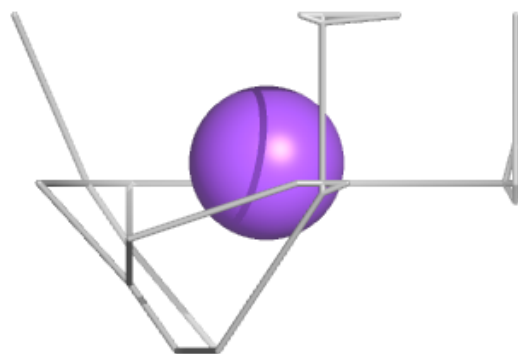
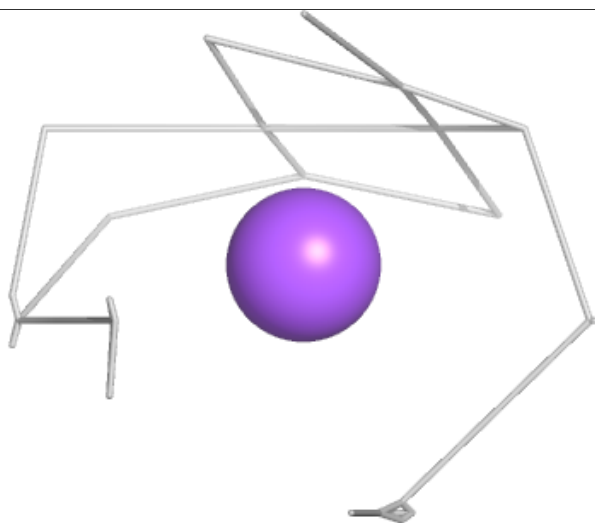
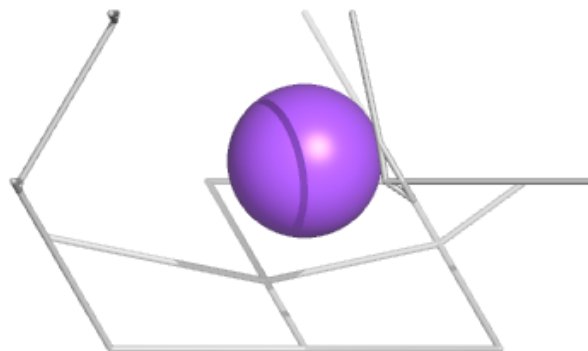
**Electron density around NA B 501:**

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and green (positive)



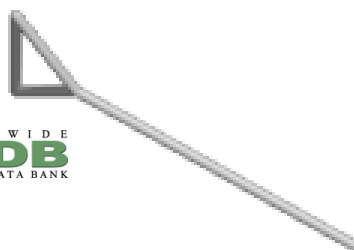
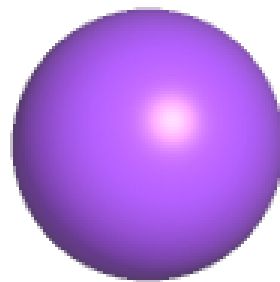
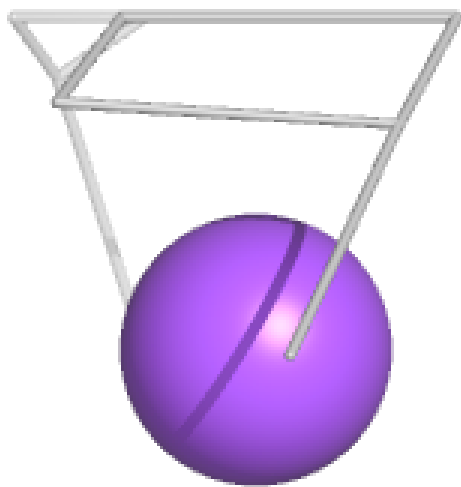
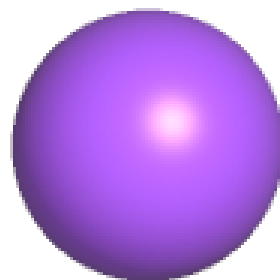
**Electron density around NA A 502:**

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and green (positive)

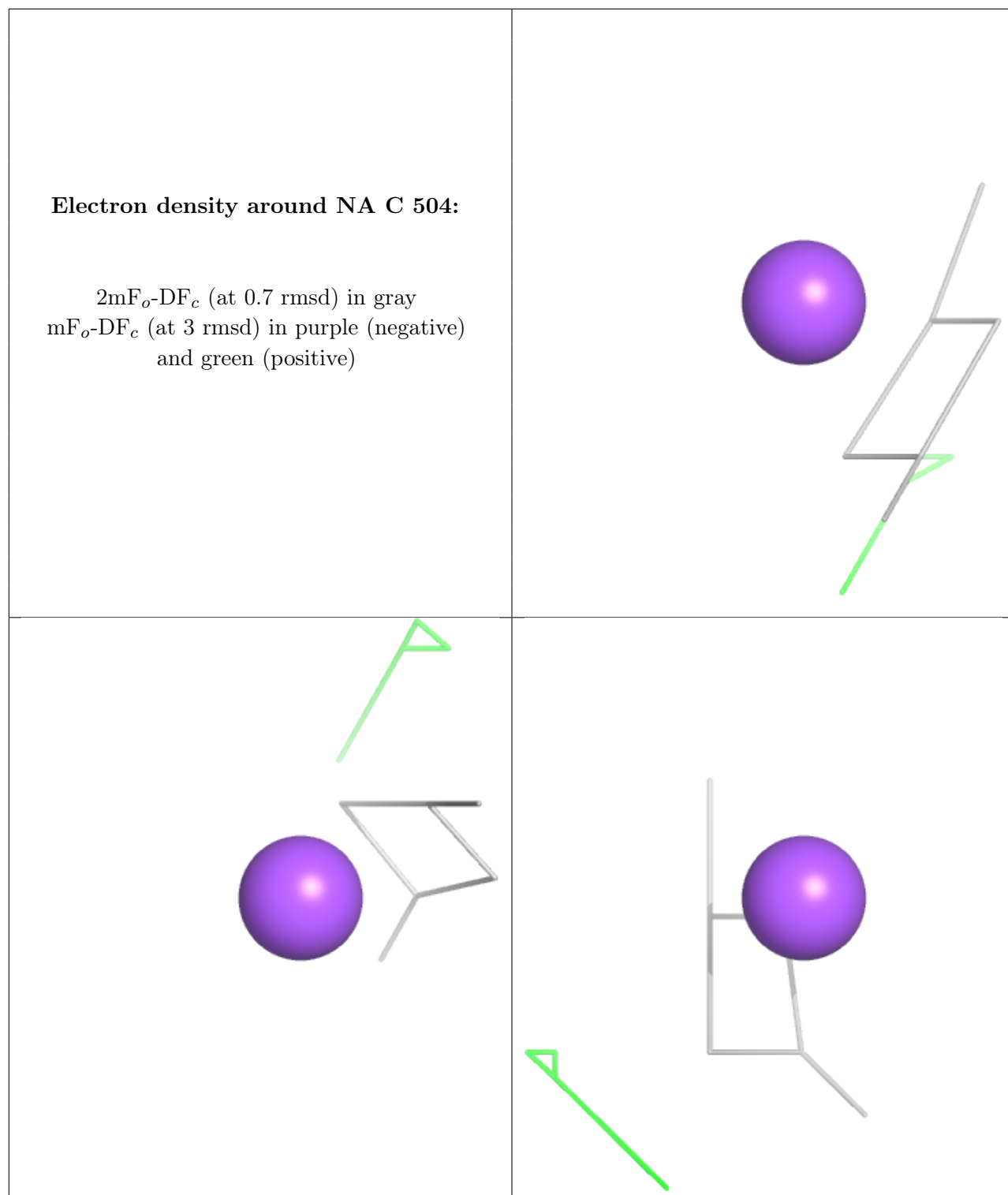


**Electron density around NA B 502:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)







## 6.5 Other polymers ⓘ

There are no such residues in this entry.